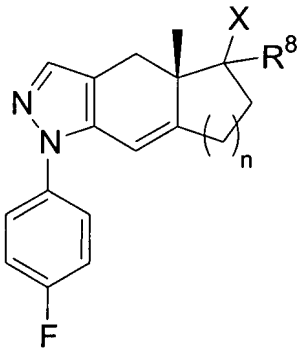
Amendments to the Claims:

This listing of claims replaces all prior versions, and listings, of claims in the application:

Listing of Claims:

1 to 7. (canceled)

8. (currently amended) A compound ~~according to Claim 1~~ of Formula II:



II

or a pharmaceutically acceptable salt or hydrate thereof, wherein:

X is selected from the group consisting of:-OR^a, -$N(R^b)$-Y-R^c, -$S(O)_j$-R^d, wherein:

Y is selected from a bond, –C(O)-, –C(O)–O–, wherein the point of attachment of the "–O–" group is to R^c forming an alkoxy moiety, -$S(O)_2$- and -C(O)-$N(R^{12})$-, wherein the point of attachment of the nitrogen group is to R^c, and

j is 0, 1 or 2,

n is 1 ~~or 2~~,

R^a, R^b, R^c, R^d and R^8 are each independently selected from the group consisting of:

(1) hydrogen, except that R^d is not hydrogen and R^c is hydrogen only when Y is a bond or -C(O)-N(R^{12})-,

(2) C_{1-6}alkyl,

(3) C_{2-6}alkenyl,

(4) C_{2-6}akynyl,

(5) C_{3-6}cycloalkyl,

(6) aryl,

(7) aralkyl,

(8) HET^1,

(9) -C_{1-6}alkyl-HET^2,

(10) aralkenyl,

(11) aralkynyl and

(12) arylsulfonylalkyl,

wherein items (2) to (5) above and the alkyl portions of items (7), (9) and (12) above and the alkenyl portion of item (10) above and the alkynyl portion of item (11) above are optionally substituted with oxo and optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{11}, $N(R^{12})_2$, C_{3-6}cycloalkyl and C_{1-4}alkyl-$S(O)_m$-, wherein m is 0, 1 or 2, and

wherein items (6) and (8) above and the aryl portion of items (7), (10), (11) and (12) above and the HET^2 portion of item (9) above are optionally substituted with one to five substituents independently selected from the group consisting of:

(a) halo,

(b) OR^{11},

(c) $N(R^{12})_2$,

(d) C_{1-6}alkyl,

(e) C_{2-6}alkenyl,

(f) C_{2-6}akynyl,

(g) C_{1-6}alkyl-$S(O)_p$-, wherein p is 0, 1 or 2,

(h) aryl,

(i) aryl-$S(O)_q$-, wherein q is 0, 1 or 2,

(j) HET^3,

(k) aralkyl,

(l) aroyl,
(m) aryloxy,
(n) aralkoxy and
(o) CN,

wherein items (d) to (g) above and the alkyl portions of item (k) above are optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{11} and $N(R^{12})_2$, and

wherein items (h), (i), (j), (l) and (m) above and the aryl portions of items (k) and (n) above are optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{12} and C_{1-4}alkyl,

each R^{11} and R^{12} is independently selected from the group consisting of hydrogen and C_{1-4}alkyl, optionally substituted with 1 to 3 halo groups; and

HET^1, HET^2 and HET^3 are each independently selected from the group of heterocycles consisting of: benzimidazolyl, benzofuranyl, benzopyrazolyl, benzotriazolyl, benzothiophenyl, benzoxazolyl, carbazolyl, carbolinyl, cinnolinyl, furanyl, imidazolyl, indolinyl, indolyl, indolazinyl, indazolyl, isobenzofuranyl, isoindolyl, isoquinolyl, isothiazolyl, isoxazolyl, naphthyridinyl, oxadiazolyl, oxazolyl, pyrazinyl, pyrazolyl, pyridopyridinyl, pyridazinyl, pyridyl, pyrimidyl, pyrrolyl, quinazolinyl, quinolyl, quinoxalinyl, thiadiazolyl, thiazolyl, thienyl, triazolyl, azetidinyl, 1,4-dioxanyl, hexahydroazepinyl, piperazinyl, piperidinyl, pyrrolidinyl, morpholinyl, thiomorpholinyl, dihydrobenzimidazolyl, dihydrobenzofuranyl, dihydrobenzothiophenyl, dihydrobenzoxazolyl, dihydrofuranyl, dihydroimidazolyl, dihydroindolyl, dihydroisooxazolyl, dihydroisothiazolyl, dihydrooxadiazolyl, dihydrooxazolyl, dihydropyrazinyl, dihydropyrazolyl, dihydropyridinyl, dihydropyrimidinyl, dihydropyrrolyl, dihydroquinolinyl, dihydrotetrazolyl, dihydrothiadiazolyl, dihydrothiazolyl, dihydrothienyl, dihydrotriazolyl, dihydroazetidinyl, methylenedioxybenzoyl, tetrahydrofuranyl, and tetrahydrothienyl.

9. (currently amended) The compound according to Claim 8 wherein

X is $-OR^a$

~~n is 1~~, and

R^a is selected from the group consisting of:

(1) hydrogen,
(2) acetyl,
(3) benzyl,
(4) C_{1-6}alkyl,
(5) C_{2-6}alkenyl,
(6) C_{2-6}alkynyl and
(7) C_{3-6}cycloalkyl,

R^8 is selected from the group consisting of:

(1) hydrogen,
(2) C_{1-6}alkyl,
(3) C_{2-6}alkenyl,
(4) C_{2-6}akynyl,
(5) C_{3-6}cycloalkyl,
(6) aryl,
(7) aralkyl,
(8) HET^1,
(9) -C_{1-6}alkyl-HET^2,
(10) aralkenyl,
(11) aralkynyl, and
(12) arylsulfonylalkyl

wherein items (2) to (5) above and the alkyl portions of items (7), (9) and (12) above and the alkenyl portion of item (10) above and the alkynyl portion of item (11) above are optionally substituted with oxo and optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{11} and C_{3-6}cycloalkyl,

wherein items (6) and (8) above and aryl portion of items (7), (10), (11) and (12) above and the HET^2 portion of item (9) above are optionally substituted with one to five substituents independently selected from the group consisting of:

(a) halo,
(b) C_{1-6}alkyl,
(c) C1-4alkoxy and
(d) aryl,

R^{11} is selected from the group consisting of hydrogen and C_{1-4}alkyl, optionally substituted with 1 to 3 halo groups; and

HET^1 and HET^2 are each independently selected from the group of heterocycles consisting of: benzimidazolyl, benzofuranyl, benzopyrazolyl, benzotriazolyl, benzothiophenyl, benzoxazolyl, carbazolyl, carbolinyl, cinnolinyl, furanyl, imidazolyl, indolinyl, indolyl, indolazinyl, indazolyl, isobenzofuranyl, isoindolyl, isoquinolyl, isothiazolyl, isoxazolyl, naphthyridinyl, oxadiazolyl, oxazolyl, pyrazinyl, pyrazolyl, pyridopyridinyl, pyridazinyl, pyridyl, pyrimidyl, pyrrolyl, quinazolinyl, quinolyl, quinoxalinyl, thiadiazolyl, thiazolyl, thienyl, triazolyl, azetidinyl, 1,4-dioxanyl, hexahydroazepinyl, piperazinyl, piperidinyl, pyrrolidinyl, morpholinyl, thiomorpholinyl, dihydrobenzimidazolyl, dihydrobenzofuranyl, dihydrobenzothiophenyl, dihydrobenzoxazolyl, dihydrofuranyl, dihydroimidazolyl, dihydroindolyl, dihydroisooxazolyl, dihydroisothiazolyl, dihydrooxadiazolyl, dihydrooxazolyl, dihydropyrazinyl, dihydropyrazolyl, dihydropyridinyl, dihydropyrimidinyl, dihydropyrrolyl, dihydroquinolinyl, dihydrotetrazolyl, dihydrothiadiazolyl, dihydrothiazolyl, dihydrothienyl, dihydrotriazolyl, dihydroazetidinyl, methylenedioxybenzoyl, tetrahydrofuranyl, and tetrahydrothienyl.

10. (original) The compound according to Claim 9 wherein:

R^8 is selected from the group consisting of:

(1) hydrogen,
(2) C_{1-6}alkyl,
(3) C_{2-6}alkenyl,
(4) C_{2-6}akynyl,
(5) C_{3-6}cycloalkyl,
(6) phenyl or naphthyl,
(7) benzyl or phenethyl,
(8) benzothiophene,

(9) phenylethenyl,
(10) phenylethynyl, and
(11) phenylsulfonylmethyl,

wherein items (2) to (5) above are optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{11} and C_{3-6}cycloalkyl,

wherein item (6) above and the phenyl portions of items (7), (9), (10) and (11) above are optionally substituted with one to five substituents independently selected from the group consisting of:

(a) halo,
(b) C_{1-6}alkyl,
(c) C_{1-4}alkoxy and
(d) phenyl.

11. (currently amended) The compound according to Claim 8 wherein:

X is $-N(R^b)-Y-R^c$, wherein:

Y is selected from –C(O)-, –C(O)–O–, wherein the point of attachment of the "–O–" group is to R^c forming an alkoxy moiety, $-S(O)_2-$ and $-C(O)-N(R^{12})-$, wherein the point of attachment of the nitrogen group is to R^c, ~~and~~

~~n is 1,~~

R^8 is hydrogen, and

R^b and R^c are each independently selected from the group consisting of:

(1) hydrogen, except that R^c is not hydrogen,
(2) C_{1-6}alkyl,
(3) C_{2-6}alkenyl,
(4) C_{2-6}akynyl,
(5) C_{3-6}cycloalkyl,

(6) aryl,
(7) aralkyl,
(8) HET^1,
(9) $-C_{1-6}$alkyl-HET^2,
(10) aralkenyl,
(11) aralkynyl and
(12) arylsulfonylalkyl,

wherein items (2) to (5) above and the alkyl portions of items (7), (9) and (12) above and the alkenyl portion of item (10) above and the alkynyl portion of item (11) above are optionally substituted with oxo and optionally substituted with with one to three substituents independently selected from the group consisting of: halo, OR^{11}, $N(R^{12})_2$, C_{3-6}cycloalkyl and C_{1-4}alkyl-$S(O)_m$-, wherein m is 0, 1 or 2, and

wherein items (6) and (8) above and the aryl portion of items (7), (10), (11) and (12) above and the HET^2 portion of item (9) above are optionally substituted with one to five substituents independently selected from the group consisting of:

(a) halo,
(b) OR^{11},
(c) $N(R^{12})_2$,
(d) C_{1-6}alkyl,
(e) C_{2-6}alkenyl,
(f) C_{2-6}akynyl,
(g) C_{1-6}alkyl-$S(O)_p$-, wherein p is 0, 1 or 2,
(h) aryl,
(i) aryl-$S(O)_q$-, wherein q is 0, 1 or 2,
(j) HET^3,
(k) aralkyl,
(l) aroyl,
(m) aryloxy,
(n) aralkoxy and
(o) CN,

wherein items (d) to (g) above and the alkyl portions of item (k) above are optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{11} and $N(R^{12})_2$, and

wherein items (h), (i), (j), (l) and (m) above and the aryl portions of items (k) and (n) above are optionally substituted with one to three substituents independently selected from the group consisting of: halo, OR^{12} and C_{1-4}alkyl,

each R^{11} and R^{12} is independently selected from the group consisting of hydrogen and C_{1-4}alkyl, optionally substituted with 1 to 3 halo groups; and

HET^1, HET^2 and HET^3 are each independently selected from the group of heterocycles consisting of: benzimidazolyl, benzofuranyl, benzopyrazolyl, benzotriazolyl, benzothiophenyl, benzoxazolyl, carbazolyl, carbolinyl, cinnolinyl, furanyl, imidazolyl, indolinyl, indolyl, indolazinyl, indazolyl, isobenzofuranyl, isoindolyl, isoquinolyl, isothiazolyl, isoxazolyl, naphthyridinyl, oxadiazolyl, oxazolyl, pyrazinyl, pyrazolyl, pyridopyridinyl, pyridazinyl, pyridyl, pyrimidyl, pyrrolyl, quinazolinyl, quinolyl, quinoxalinyl, thiadiazolyl, thiazolyl, thienyl, triazolyl, azetidinyl, 1,4-dioxanyl, hexahydroazepinyl, piperazinyl, piperidinyl, pyrrolidinyl, morpholinyl, thiomorpholinyl, dihydrobenzimidazolyl, dihydrobenzofuranyl, dihydrobenzothiophenyl, dihydrobenzoxazolyl, dihydrofuranyl, dihydroimidazolyl, dihydroindolyl, dihydroisooxazolyl, dihydroisothiazolyl, dihydrooxadiazolyl, dihydrooxazolyl, dihydropyrazinyl, dihydropyrazolyl, dihydropyridinyl, dihydropyrimidinyl, dihydropyrrolyl, dihydroquinolinyl, dihydrotetrazolyl, dihydrothiadiazolyl, dihydrothiazolyl, dihydrothienyl, dihydrotriazolyl, dihydroazetidinyl, methylenedioxybenzoyl, tetrahydrofuranyl, and tetrahydrothienyl.

12. (original) The compound according to Claim 11 wherein:

R^b and R^c are each independently selected from the group consisting of:

(1) hydrogen, except that R^c is hydrogen only when Y is a bond or -C(O)-$N(R^{12})$-,

(2) C_{1-6}alkyl,
(3) C_{2-6}alkenyl,
(4) C_{2-6}akynyl,
(5) C_{3-6}cycloalkyl,
(6) aryl,
(7) aralkyl,
(8) HET^1,
(9) -C_{1-6}alkyl-HET^2,
(10) aralkenyl, and
(11) aralkynyl,

wherein items (2) to (5) above are optionally substituted with 1-3 halo groups, and

wherein items (6) and (8) and aryl portion of items (7), (10) and (11) above and the HET^2 portion of item (9) above are optionally substituted with one to five substituents independently selected from the group consisting of:

(a) halo,
(b) C_{1-4}alkyl, optionally substituted with 1-3 halo groups, and
(c) C_{1-4}alkylthio,

HET^1 and HET^2 are each independently selected from the group of heterocycles consisting of: benzimidazolyl, benzofuranyl, benzopyrazolyl, benzotriazolyl, benzothiophenyl, benzoxazolyl, carbazolyl, carbolinyl, cinnolinyl, furanyl, imidazolyl, indolinyl, indolyl, indolazinyl, indazolyl, isobenzofuranyl, isoindolyl, isoquinolyl, isothiazolyl, isoxazolyl, naphthyridinyl, oxadiazolyl, oxazolyl, pyrazinyl, pyrazolyl, pyridopyridinyl, pyridazinyl, pyridyl, pyrimidyl, pyrrolyl, quinazolinyl, quinolyl, quinoxalinyl, thiadiazolyl, thiazolyl, thienyl, triazolyl, azetidinyl, 1,4-dioxanyl, hexahydroazepinyl, piperazinyl, piperidinyl, pyrrolidinyl, morpholinyl, thiomorpholinyl, dihydrobenzimidazolyl, dihydrobenzofuranyl, dihydrobenzothiophenyl, dihydrobenzoxazolyl, dihydrofuranyl, dihydroimidazolyl, dihydroindolyl, dihydroisooxazolyl, dihydroisothiazolyl, dihydrooxadiazolyl, dihydrooxazolyl, dihydropyrazinyl, dihydropyrazolyl, dihydropyridinyl, dihydropyrimidinyl, dihydropyrrolyl, dihydroquinolinyl, dihydrotetrazolyl, dihydrothiadiazolyl, dihydrothiazolyl, dihydrothienyl, dihydrotriazolyl, dihydroazetidinyl, methylenedioxybenzoyl, tetrahydrofuranyl, and tetrahydrothienyl.

13. (original) The compound according to Claim 12 wherein:

R^b is selected from the group consisting of: hydrogen and C_{1-4}alkyl, and

R^c is selected from the group consisting of:

(1) C_{1-4}alkyl,

(2) phenyl or benzyl, each optionally substituted with 1 to 5 groups independently selected from fluoro, chloro and trifluoromethyl,

(3) naphthyl,

(4) thiopheneyl,

(5) pyridyl,

(6) isoquinolyl,

(7) pyrimidyl and

(8) pyrazyl,

wherein items (4) to (8) above are optionally substitited with 1 to 5 groups independently selected from fluoro, chloro, methyl,methylthio and trifluoromethyl.

14. (original) The compound according to Claim 13, wherein R^c is phenyl, optionally substituted with 1 to 5 groups independently selected from fluoro, chloro and trifluoromethyl.

15. (currently amended) The compound according to Claim 8, wherein:

X is $-S(O)_j-R^d$, wherein j is 0, 1 or 2,

~~n is 1,~~

R^8 is hydrogen, and

R^d is selected from the group consisting of:

(1) C_{1-6}alkyl,

(2) C_{2-6}alkenyl,
(3) C_{2-6}akynyl,
(4) C_{3-6}cycloalkyl,
(5) aryl,
(6) aralkyl,
(7) HET^1,
(8) -C_{1-6}alkyl-HET^2,
(9) aralkenyl, and
(10) aralkynyl,

wherein items (1) to (4) above are optionally substituted with 1-3 halo groups, and

wherein items (5) and (7) and aryl portion of items (6), (9) and (10) above and the HET^2 portion of item (8) above are optionally substituted with one to five substituents independently selected from the group consisting of:

(a) halo,
(b) C_{1-4}alkyl, optionally substituted with 1-3 halo groups, and
(c) C_{1-4}alkylthio, and

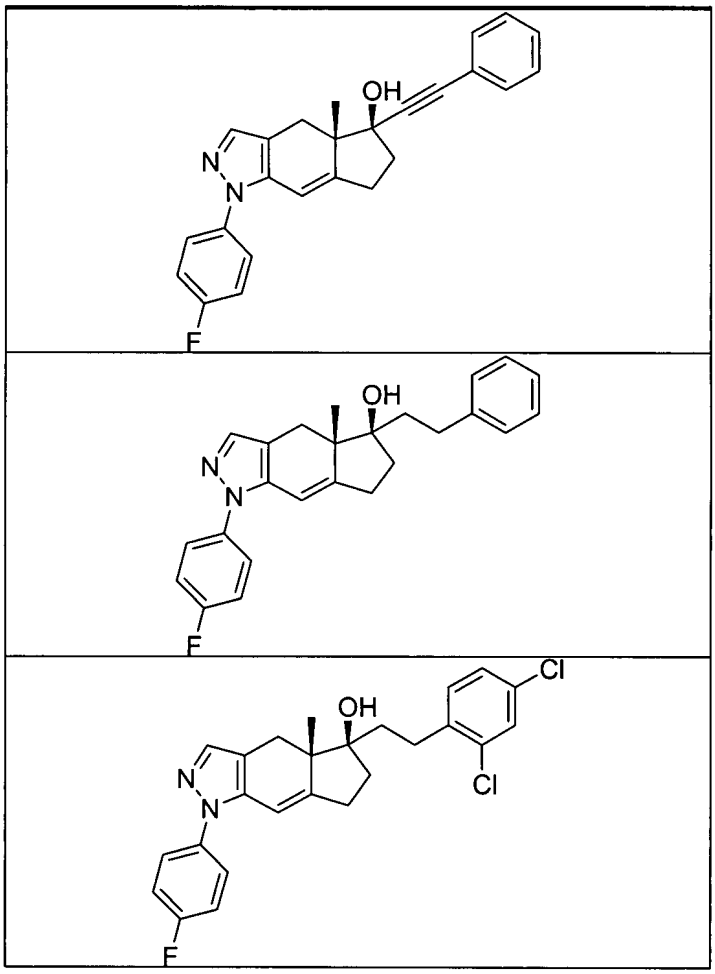
HET^1 and HET^2 are each independently selected from the group of heterocycles consisting of: benzimidazolyl, benzofuranyl, benzopyrazolyl, benzotriazolyl, benzothiophenyl, benzoxazolyl, carbazolyl, carbolinyl, cinnolinyl, furanyl, imidazolyl, indolinyl, indolyl, indolazinyl, indazolyl, isobenzofuranyl, isoindolyl, isoquinolyl, isothiazolyl, isoxazolyl, naphthyridinyl, oxadiazolyl, oxazolyl, pyrazinyl, pyrazolyl, pyridopyridinyl, pyridazinyl, pyridyl, pyrimidyl, pyrrolyl, quinazolinyl, quinolyl, quinoxalinyl, thiadiazolyl, thiazolyl, thienyl, triazolyl, azetidinyl, 1,4-dioxanyl, hexahydroazepinyl, piperazinyl, piperidinyl, pyrrolidinyl, morpholinyl, thiomorpholinyl, dihydrobenzimidazolyl, dihydrobenzofuranyl, dihydrobenzothiophenyl, dihydrobenzoxazolyl, dihydrofuranyl, dihydroimidazolyl, dihydroindolyl, dihydroisooxazolyl, dihydroisothiazolyl, dihydrooxadiazolyl, dihydrooxazolyl, dihydropyrazinyl, dihydropyrazolyl, dihydropyridinyl, dihydropyrimidinyl, dihydropyrrolyl, dihydroquinolinyl, dihydrotetrazolyl, dihydrothiadiazolyl, dihydrothiazolyl, dihydrothienyl, dihydrotriazolyl, dihydroazetidinyl, methylenedioxybenzoyl, tetrahydrofuranyl, and tetrahydrothienyl.

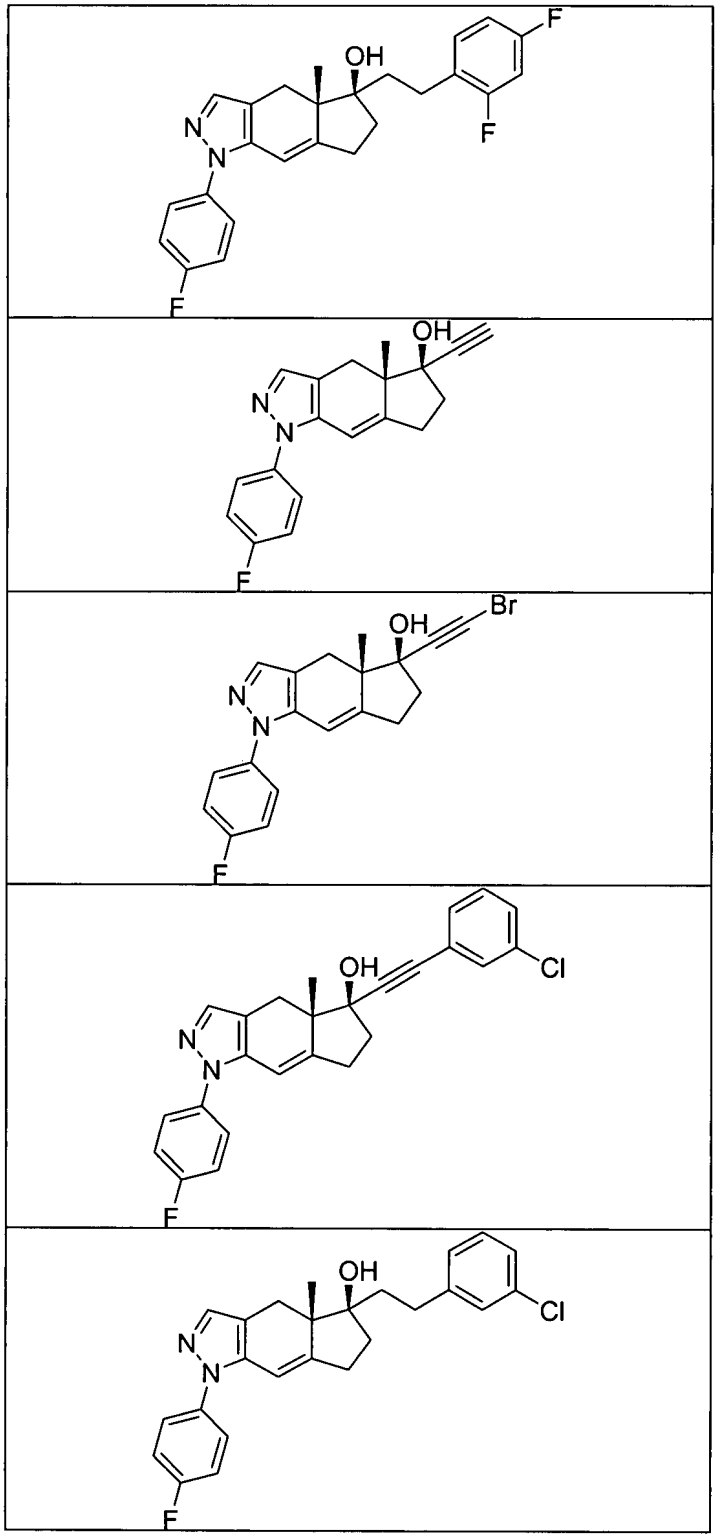
16. (original) The compound according to Claim 15, wherein

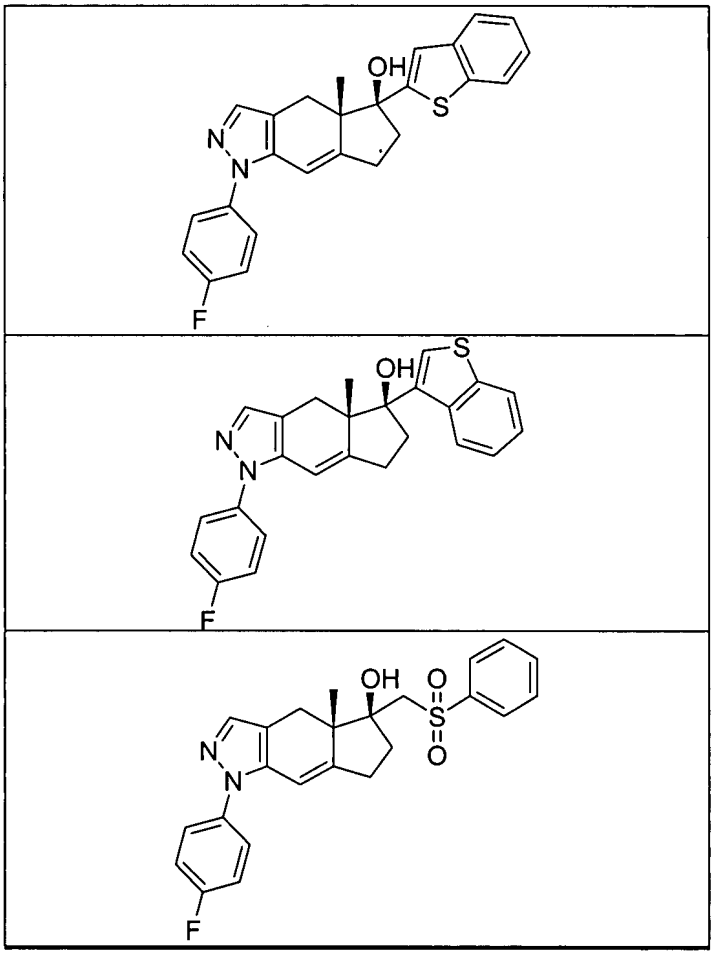
R^d is phenyl, optionally substituted with 1 to 5 groups independently selected from fluoro, chloro and trifluoromethyl.

17. (currently amended) A compound selected from the following group:

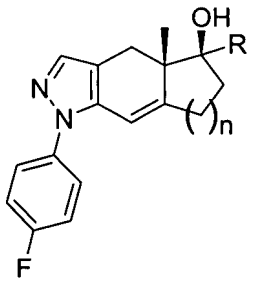
~~Table A~~




OH
F
F
N
N
F
OH
N
N
F
OH
Br
N
N
F
OH
Cl
N
N
F
OH
Cl
N
N
F



~~Table B~~

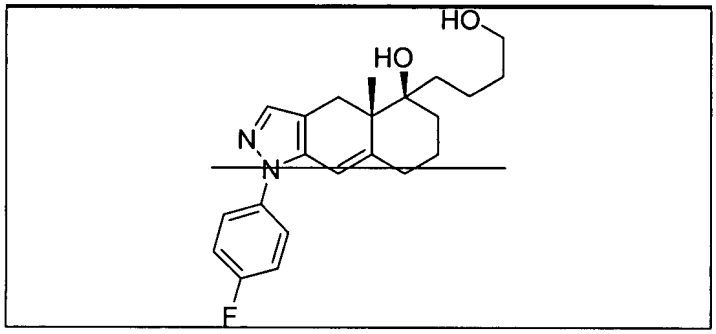


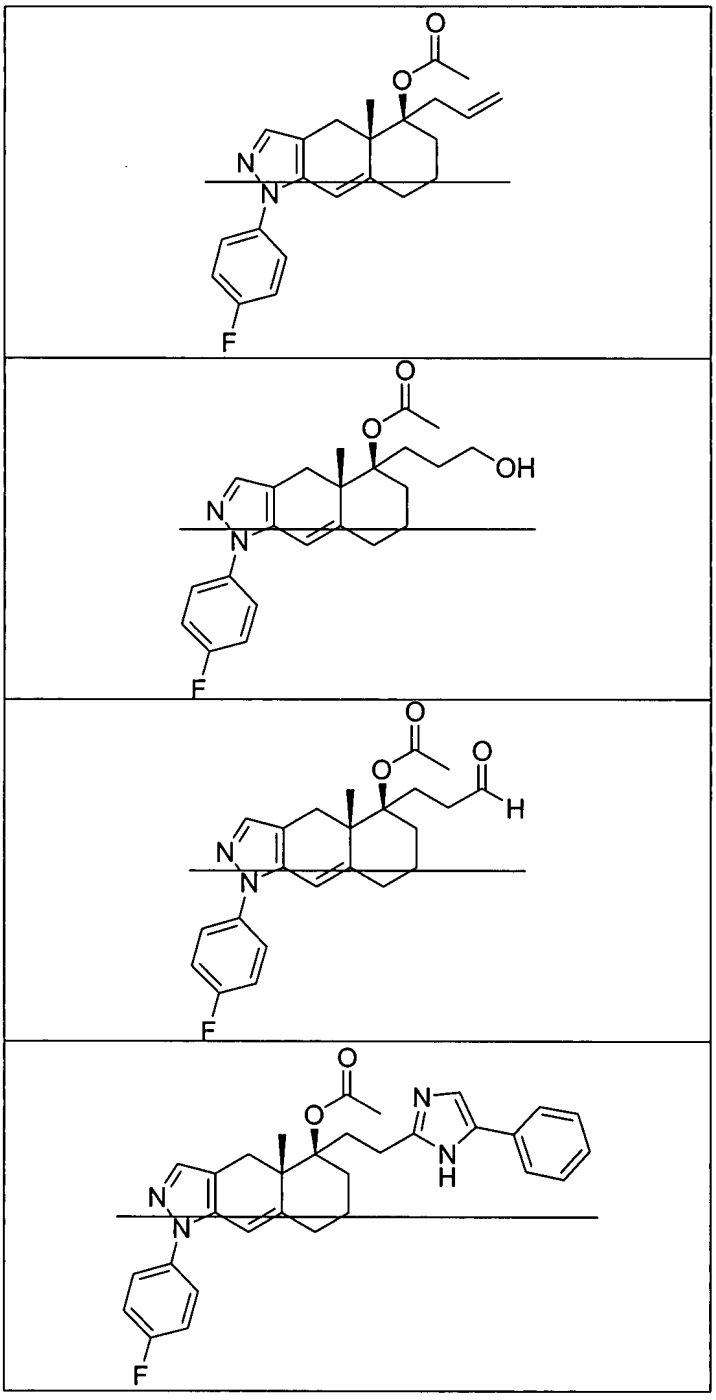
n	R Group
1	Vinyl
~~2~~	~~Vinyl~~
1	Allyl

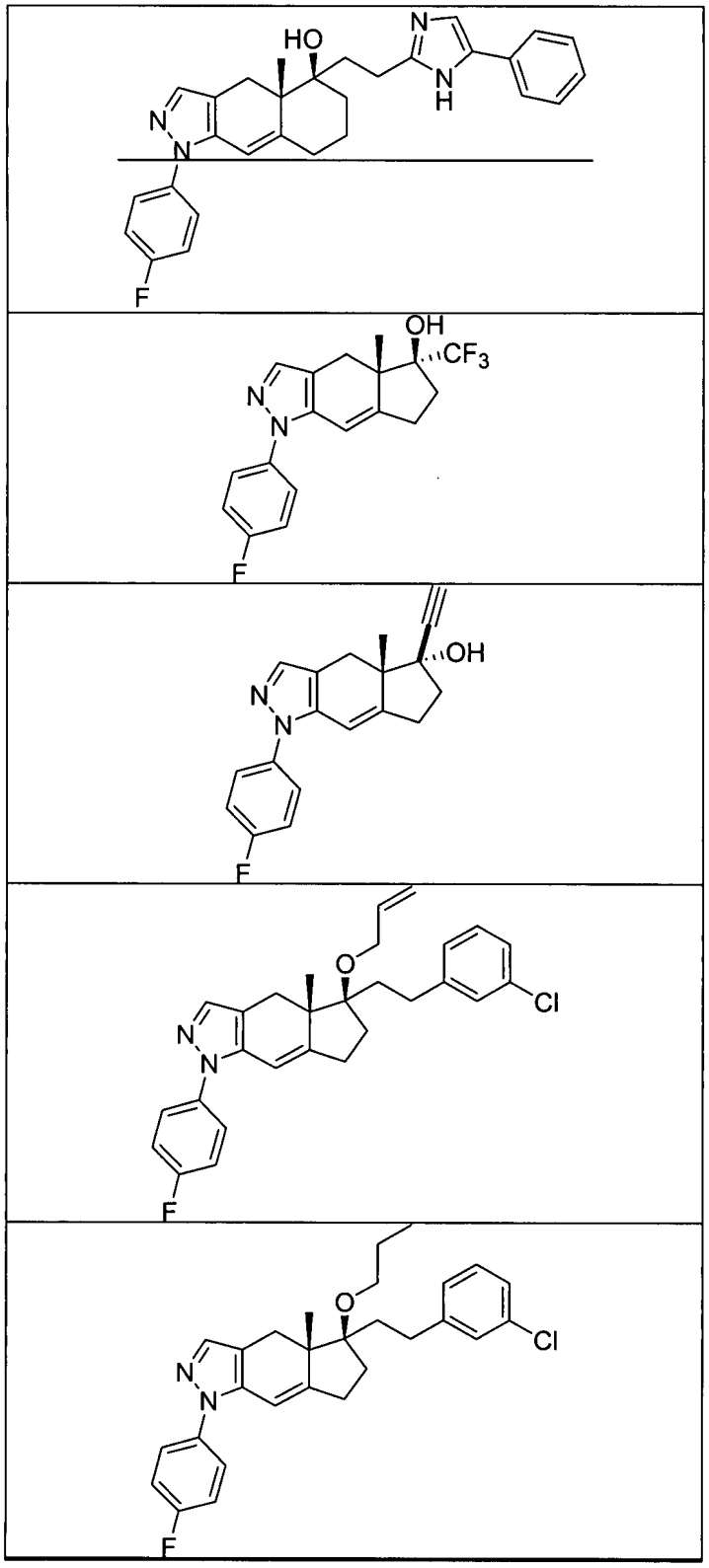
~~2~~	~~Allyl~~
1	3-butenyl
~~2~~	~~3-butenyl~~
~~2~~	~~*n*-butyl~~
1	*n*-pentyl
~~2~~	~~*n*-pentyl~~
~~2~~	~~*n*-hexyl~~
1	3-methylbutyl
~~2~~	~~3-methylbutyl~~
1	2-cyclohexylethyl
~~2~~	~~2-cyclohexylethyl~~
1	3,3-dimethylbutyl
~~2~~	~~3,3-dimethylbutyl~~
1	4-methyl-3-pentenyl
~~2~~	~~4-methyl-3-pentenyl~~
1	4,4,4-trifluorobutyl
~~2~~	~~4,4,4-trifluorobutyl~~
1	3,4,4-trifluoro-3-butenyl
~~2~~	~~3,4,4-trifluoro-3-butenyl~~
1	3-methoxypropyl
~~2~~	~~3-methoxypropyl~~
~~2~~	~~benzyl~~
~~2~~	~~phenyl~~
~~2~~	~~phenethyl~~
~~2~~	~~3-phenylpropyl~~
1	2-(2-chlorophenyl)ethyl
~~2~~	~~2-(2-chlorophenyl)ethyl~~
~~2~~	~~2-(3-chlorophenyl)ethyl~~
1	2-(4-chlorophenyl)ethyl
~~2~~	~~2-(4-chlorophenyl)ethyl~~
~~2~~	~~2-(2,4-dichlorophenyl)ethyl~~
1	2-(4-fluorophenyl)ethyl
2	2-(4-fluorophenyl)ethyl
1	2-(2,5-difluorophenyl)ethyl
1	2-(2,3-difluorophenyl)ethyl
1	2-(3,5-difluorophenyl)ethyl
1	2-(4-methoxyphenyl)ethyl
~~2~~	~~2-(4-methoxyphenyl)ethyl~~
1	2-(2-naphthyl)ethyl
~~2~~	~~2-(2-naphthyl)ethyl~~
~~2~~	~~2-(2,4-difluorophenyl)ethyl~~
1	2-(3-(trifluoromethyl)phenyl)ethyl

~~2~~	~~2-(3-(trifluoromethyl)phenyl)ethyl~~
1	2-(2-methoxyphenyl)ethyl
~~2~~	~~2-(2-methoxyphenyl)ethyl~~
1	2-(4-*tert*-butylphenyl)ethyl
~~2~~	~~2-(4-*tert*-butylphenyl)ethyl~~
1	2-(4-methylphenyl)ethyl
~~2~~	~~2-(4-methylphenyl)ethyl~~
1	2-(1-naphthyl)ethyl
~~2~~	~~2-(1-naphthyl)ethyl~~
1	2-(2-methylphenyl)ethyl
~~2~~	~~2-(2-methylphenyl)ethyl~~
1	2-(3-methylphenyl)ethyl
~~2~~	~~2-(3-methylphenyl)ethyl~~
1	2-(2-fluorophenyl)ethyl
1	2-(3-fluorophenyl)ethyl
1	2-(3,4-dichlorophenyl)ethyl
1	2-(2-chloro-4-fluorophenyl)ethyl
1	2-(3-thiopheneyl)ethyl
1	3-(*N*-pyrrolyl)propyl
~~2~~	~~3-(*N*-pyrrolyl)propyl~~
1	*E*-2-phenylethenyl
~~2~~	~~*E*-2-phenylethenyl~~
1	*Z*-2-phenylethenyl
~~2~~	~~*Z*-2-phenylethenyl~~
~~2~~	~~2-phenylethynyl~~
1	2-(2,4-difluorophenyl)ethynyl
1	2-(2-thiopheneyl)ethyl
1	2-(3,4-difluorophenyl)ethyl
1	2-(3,4,5-trifluorophenyl)ethyl
~~2~~	~~H~~
1	H

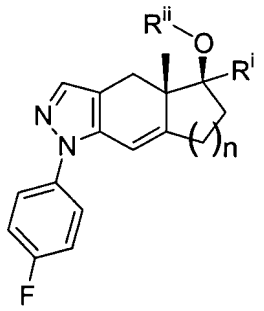
~~Table C~~




O
O
N
N
F
O
O
OH
N
N
F
O
O
O
H
N
N
F
O
O
N
N
H
N
N
F


HO
N
N
H
N
N
F
OH
CF3
N
N
F
OH
N
N
F
O
Cl
N
N
F
O
Cl
N
N
F

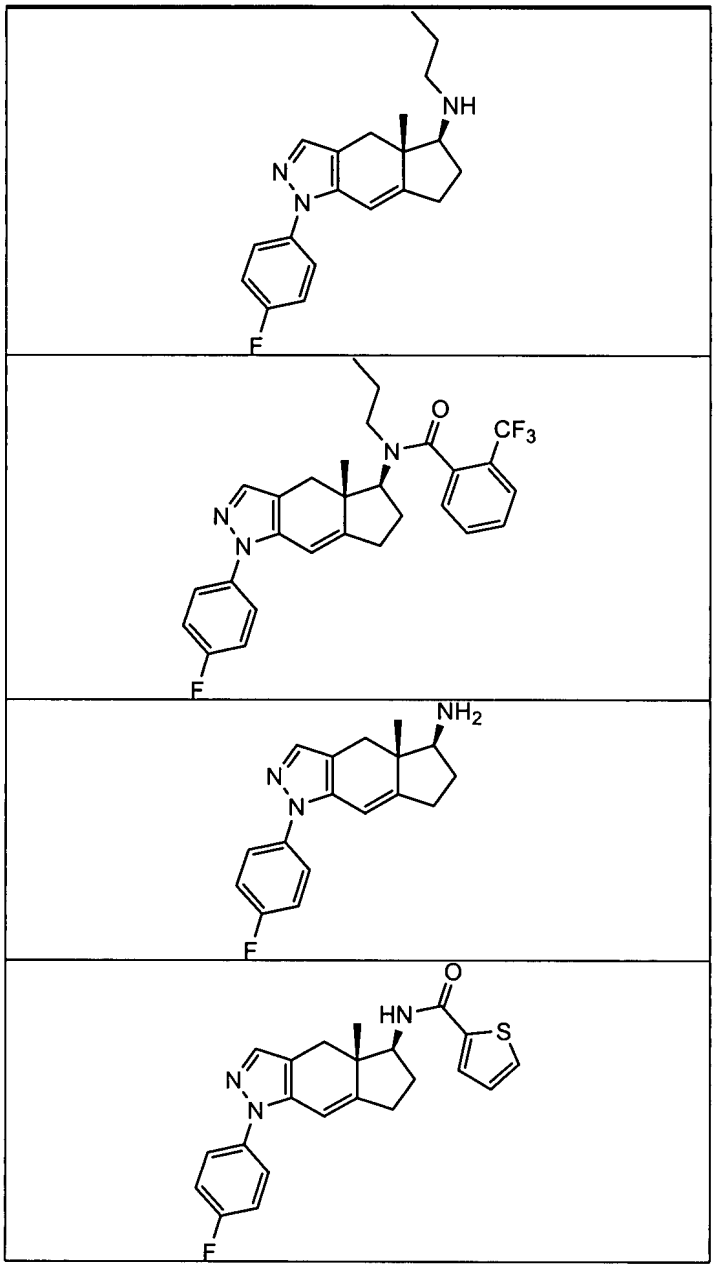
~~Table D~~

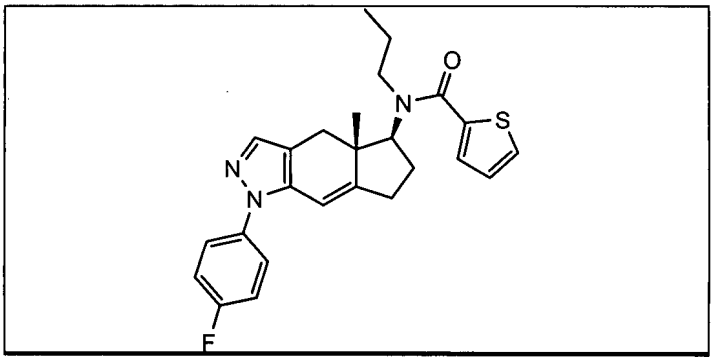


n	R^i Group	R^{ii} Group
1	H	*n*-propyl
~~2~~	~~allyl~~	~~allyl~~
1	phenethyl	cyclopropylmethyl
1	phenethyl	*E*-2-butenyl
1	2-(2,4-dichlorophenyl)ethyl	methyl
1	2-(2-chlorophenyl)ethyl	methyl
1	2-(3-chlorophenyl)ethyl	methyl
1	*n*-pentyl	methyl
1	2-(4-fluorophenyl)ethyl	methyl
~~2~~	~~phenethyl~~	~~methyl~~
1	2-(2,4-dichlorophenyl)ethyl	benzyl
1	2-(2,4-dichlorophenyl)ethyl	allyl
1	2-(2,4-dichlorophenyl)ethyl	*n*-propyl
1	2-(2-chlorophenyl)ethyl	*n*-propyl
1	phenethyl	*n*-propyl
1	phenethyl	methyl
1	2-(3-chlorophenyl)ethynyl	allyl
1	2-(3-chlorophenyl)ethynyl	*n*-propyl
1	2-(2,4-difluorophenyl)ethyl	methyl
~~2~~	~~2-(2,4-difluorophenyl)ethyl~~	~~methyl~~
1	phenethyl	*E*-2-pentenyl
1	trifluoromethyl	allyl
1	trifluoromethyl	*n*-propyl
1	2-(3-methylphenyl)ethyl	methyl
1	phenethyl	*n*-butyl
1	phenethyl	*n*-pentyl
1	2-(3,4-difluorophenyl)ethyl	*n*-propyl

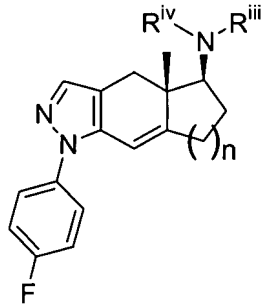
1	2-(3-fluorophenyl)ethyl	methyl

~~Table E~~





~~Table F~~

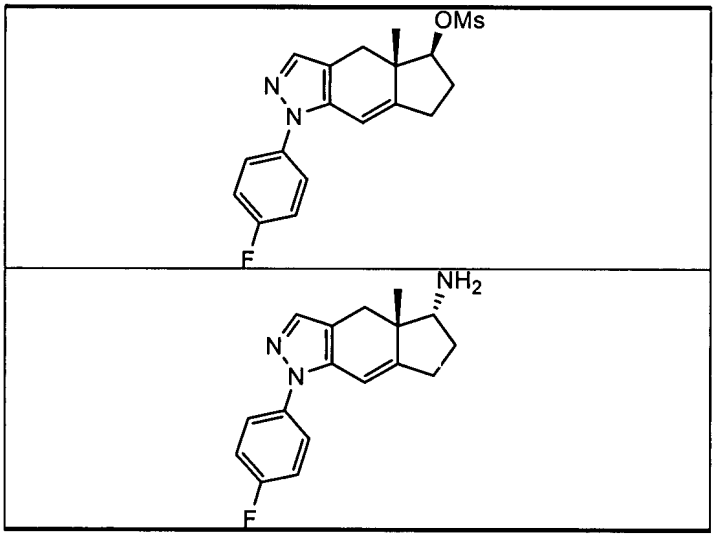


n	R^{iii} Group	R^{iv} Group
1	CO(3,3,3-trifluoropropyl)	H
1	CO(2-(trifluoromethyl)phenyl)	H
1	CO(3-chlorophenyl)	H
1	$COCH_2$(2-(trifluoromethyl)phenyl)	H
1	CO(2,4,5-trifluorophenyl)	H
1	CO(3-thiopheneyl)	H
1	$COCH_2$(2,4-difluorophenyl)	H
1	$COCH_2$(3-chlorophenyl)	H
1	CO(3-chloro-4-fluorophenyl)	H
1	$COCH_2$(2,5-difluorophenyl)	H
1	$COCH_2$(2-thiopheneyl)	H
1	$COCH_2$(3-thiopheneyl)	H
1	CO(2-chloro-5-fluorophenyl)	H
1	$COCH_2$(3-chloro-4-fluorophenyl)	H
1	CO(2,4,5-trifluorophenyl)	methyl
1	CO(2-(trifluoromethyl)phenyl)	methyl
1	CO(2-thiopheneyl)	methyl

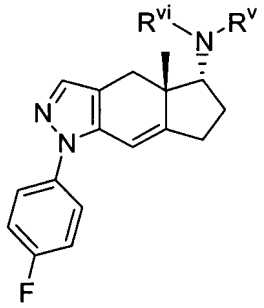
1	CO(3-chlorophenyl)	methyl
1	CO(phenyl)	H
1	CO(2,4-difluorophenyl)	H
1	$COCH_2$(3-chloro-4-fluorophenyl)	methyl
1	CO(2,4-difluorophenyl)	methyl
1	$COCH_2$(2-(trifluoromethyl)phenyl)	methyl
1	CO(2-fluorophenyl)	H
1	CO(2,6-difluorophenyl)	H
1	CO(2-chlorophenyl)	H
1	CO(1-naphthyl)	H
1	CO(2-(trifluoromethyl)-4-fluorophenyl)	H
1	CO(2,5-difluorophenyl)	H
1	CO(2,3-difluorophenyl)	H
1	CO(2-chloro-4-fluorophenyl)	H
1	CO(2-chloro-3-fluorophenyl)	H
1	CO(tert-butyl)	H
1	CO(isopropyl)	H
1	CO(2-chloro-3-fluorophenyl)	methyl
1	CO(2-(trifluoromethyl)-4-fluorophenyl)	methyl
1	CO(2,6-difluorophenyl)	methyl
1	CO(2-chloro-4-fluorophenyl)	methyl
1	SO_2(phenyl)	H
1	CO(2,6-dichlorophenyl)	methyl
1	CO(2,6-dichlorophenyl)	H
~~2~~	~~CO(phenyl)~~	~~H~~
~~2~~	~~CO(2-(trifluoromethyl)phenyl)~~	~~H~~
~~2~~	~~CO(2-chloro-4-fluorophenyl)~~	~~H~~
~~2~~	~~CO(2-chlorophenyl)~~	~~H~~
~~2~~	~~CO(2-fluorophenyl)~~	~~H~~
~~2~~	~~$COCH_2$(2-(trifluoromethyl)phenyl)~~	~~H~~
~~2~~	~~$COCH_2$(2,4-difluorophenyl)~~	~~H~~
~~2~~	~~$COCH_2$(3-chlorophenyl)~~	~~H~~
1	SO_2(2,4-difluorophenyl)	H
1	CO(2,4-difluorophenyl)	*n*-propyl
1	SO_2(3-chlorophenyl)	H
2	CO(3-chlorophenyl)	H
1	SO_2(2-chloro-4-fluorophenyl)	H
1	CO_2(phenyl)	methyl
1	CO_2(phenyl)	H
1	CONH(phenyl)	H
1	SO_2(2-fluorophenyl)	H
1	SO_2(2-chlorophenyl)	H

1	CONH(phenyl)	methyl
1	SO_2(2-(trifluoromethyl)phenyl)	H
~~2~~	~~CONH(phenyl)~~	~~H~~
1	SO_2(3-fluorophenyl)	H
~~2~~	~~CO_2(phenyl)~~	~~H~~
~~2~~	~~CO(2,4-difluorophenyl)~~	~~H~~
1	CO(2-chloro-4-fluorophenyl)	*n*-propyl
~~2~~	~~CO(2-(trifluoromethyl)-4-fluorophenyl)~~	~~H~~
~~2~~	~~CO(2-chloro-4-fluorophenyl)~~	~~methyl~~
~~2~~	~~CO(2-(trifluoromethyl)-4-fluorophenyl)~~	~~methyl~~
~~2~~	~~CO(2,4-difluorophenyl)~~	~~methyl~~
1	CO(2-(trifluoromethyl)-4-fluorophenyl)	*n*-propyl
~~2~~	~~SO_2(2-chloro-4-fluorophenyl)~~	~~H~~
~~2~~	~~SO_2(2,4-difluorophenyl)~~	~~H~~
1	SO_2(3-chlorophenyl)	methyl
1	SO_2(2-chloro-4-fluorophenyl)	methyl
1	SO_2(3-chlorophenyl)	*n*-propyl
~~2~~	~~CO(2,4-difluorophenyl)~~	~~*n*-propyl~~
~~2~~	~~CO(2-chloro-4-fluorophenyl)~~	~~*n*-propyl~~
~~2~~	~~CO(2-(trifluoromethyl)-4-fluorophenyl)~~	~~*n*-propyl~~

~~Table G~~

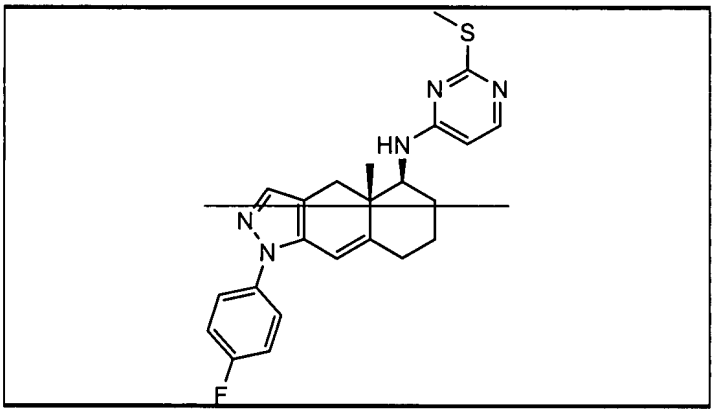


~~Table H~~

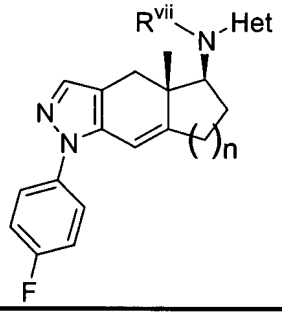


R^v Group	R^{vi} Group
CO(phenyl)	H
SO_2(2-chloro-4-fluorophenyl)	H
CO(2-chlorophenyl)	H
CO(3-chlorophenyl)	H
CO(2-(trifluoromethyl)phenyl)	H
CO(isopropyl)	H
CO(tert-butyl)	H
CO(3-thiopheneyl)	H
CO(2-thiopheneyl)	H
CO(2,4,5-trifluorophenyl)	H
CO(2,5-difluorophenyl)	H
CO_2(phenyl)	H
SO_2(phenyl)	H
CO(2-chlorophenyl)	methyl
CO(2-(trifluoromethyl)phenyl)	methyl
CO(3-chlorophenyl)	methyl
CONH(phenyl)	H
CO(2,6-difluorophenyl)	H
$COCH_2$(2,4-difluorophenyl)	H
CO(2,4-difluorophenyl)	H
CO(2-fluorophenyl)	H
CO(2-(trifluoromethyl)-4-fluorophenyl)	*n*-propyl

~~Table I~~

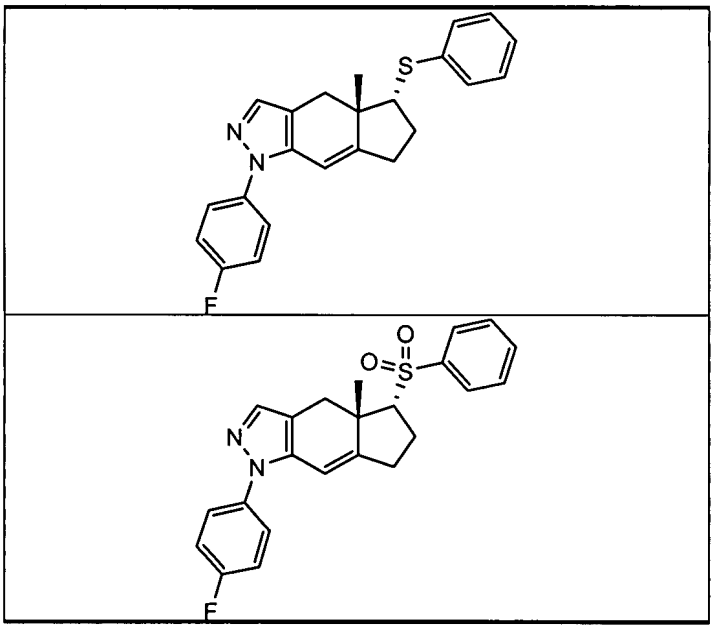


~~Table J~~

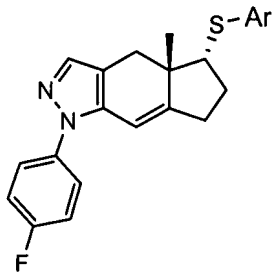


n	Het (Heterocycle)	R^{vii} Group
1	2-pyridyl	H
1	4-isoquinolyl	H
1	2-(4-methyl)pyridyl	H
1	5-pyrimidyl	H
1	3-pyridyl	H
1	2-pyrimidyl	H
1	2-pyrimidyl	methyl
1	2-pyrimidyl	*n*-propyl
1	4-(2-methylthio)pyrimidyl	H
~~2~~	~~4-(2-methylthio)pyrimidyl~~	~~methyl~~
~~2~~	~~4-(2-methylthio)pyrimidyl~~	~~*n*-propyl~~
1	2-(4-trifluoromethyl)pyridyl	H
~~2~~	~~2-(4-trifluoromethyl)pyridyl~~	~~H~~
1	2-pyrazinyl	H
1	4-(2,3,5,6-tetrafluoro)pyridyl	H
~~2~~	~~4-(2,3,5,6-tetrafluoro)pyridyl~~	~~H~~

~~Table K~~



~~Table L~~



Ar Group
2-chlorophenyl
3-chlorophenyl
2-(trifluoromethyl)phenyl
2,6-dichlorophenyl

Ar Group
2,4-dichlorophenyl

or a pharmaceutically acceptable salt of any compound selected from any of the tables above.

18. (original) A pharmaceutical composition comprising a compound according to Claim 1 in combination with a pharmaceutically acceptable carrier.

19 to 21. (canceled)